

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100220\
 Data File : VN063827.D
 Acq On : 2 Oct 2020 17:32
 Operator : JC/MD
 Sample : PB132030ZHE#06
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB132030ZHE#06

Quant Time: Oct 05 06:17:15 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 01 01:46:38 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	210518	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	403246	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	418644	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	170314	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	184678	53.10	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.20%	
35) Dibromofluoromethane	7.55	113	133756	47.80	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.60%	
50) Toluene-d8	10.06	98	515892	48.64	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.28%	
62) 4-Bromofluorobenzene	12.38	95	207319	52.00	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.00%	
Target Compounds						
16) Acetone	3.78	43	19657	18.361	ug/l	99
20) Methylene Chloride	4.51	84	20271	8.348	ug/l	98
25) 2-Butanone	6.79	43	16921	11.148	ug/l	98
43) Isopropyl Acetate	8.13	43	65836	9.834	ug/l	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN100220\
Data File : VN063827.D
Acq On : 2 Oct 2020 17:32
Operator : JC/MD
Sample : PB132030ZHE#06
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PB132030ZHE#06

Quant Time: Oct 05 06:17:15 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
Quant Title : SW846 8260
QLast Update : Thu Oct 01 01:46:38 2020
Response via : Initial Calibration

